

Report of Analysis

Client:	PSEG					Date Collected:	05/07/	/25	
Project:	OR-636 Oradell and	l New N	Ailford			Date Received:	05/08/	/25	
Client Sample ID:	OR-636-04					SDG No.:	Q1983	3	
Lab Sample ID:	Q1983-09					Matrix:	Solid		
Analytical Method:	NJEPH					% Solid:	86.8		
Sample Wt/Vol:	30.08 Units:	g				Final Vol:	2000	uL	
Soil Aliquot Vol:		uL				Test:	EPH_	NF	
Prep Method :									
Prep Date :			Date	Analyzed :				Prep Batch ID	
Prep Date : 05/09/25 09:02				Analyzed : /25 16:26				Prep Batch ID PB167926	
·				2					Datafile
05/09/25 09:02		Conc.	05/09	/25 16:26	MDL	LOQ / C	CRQL U		
05/09/25 09:02		Conc.	05/09	/25 16:26	MDL	LOQ / C	CRQL I	PB167926	
CAS Number Parameter TARGETS	hatic C28-C40	Conc. 3.31	05/09	/25 16:26	MDL 1.36	LOQ/0 2.30	CRQL U	PB167926 Units(Dry Weight)	
CAS Number Parameter TARGETS Aliphatic C28-C40 Alip			05/09	/25 16:26 Dilution		_	CRQL I	PB167926 U nits(Dry Weight) mg/kg	
CAS Number Parameter TARGETS Aliphatic C28-C40 Alip Aliphatic C9-C28 Alip	hatic C28-C40	3.31	05/09	/25 16:26 Dilution	1.36	2.30	CRQL U	PB167926 U nits(Dry Weight) mg/kg	FC068815.D

* As samples are not fractionated, all aliphatic and aromatic carbon compounds in the C9-C40 carbon range are calculated against the aliphatic calibration curve, and reported as Aliphatic EPH. Therefore, the aliphatic C9-C40 concentration for the sample is reported as the Total EPH.

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution



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Project:	OR-636 Oradell and	l New N	Ailford			Date Received:	05/08/	/25	
Client Sample ID:	OR-636-04					SDG No.:	Q1983	3	
Lab Sample ID:	Q1983-09					Matrix:	Solid		
Analytical Method:	NJEPH					% Solid:	86.8		
Sample Wt/Vol:	30.08 Units:	g				Final Vol:	2000	uL	
Soil Aliquot Vol:		uL				Test:	EPH_	NF	
Prep Method :									
Prep Date :			Date	Analyzed :				Prep Batch ID	
Prep Date : 05/09/25 09:02				Analyzed : /25 16:26				Prep Batch ID PB167926	
·				2					Datafile
05/09/25 09:02		Conc.	05/09	/25 16:26	MDL	LOQ / C	CRQL U		
05/09/25 09:02		Conc.	05/09	/25 16:26	MDL	LOQ / C	CRQL I	PB167926	
CAS Number Parameter TARGETS	hatic C28-C40	Conc. 3.31	05/09	/25 16:26	MDL 1.36	LOQ/0 2.30	CRQL U	PB167926 Units(Dry Weight)	
CAS Number Parameter TARGETS Aliphatic C28-C40 Alip			05/09	/25 16:26 Dilution		_	CRQL I	PB167926 U nits(Dry Weight) mg/kg	
CAS Number Parameter TARGETS Aliphatic C28-C40 Alip Aliphatic C9-C28 Alip	hatic C28-C40	3.31	05/09	/25 16:26 Dilution	1.36	2.30	CRQL U	PB167926 U nits(Dry Weight) mg/kg	FC068815.D

* As samples are not fractionated, all aliphatic and aromatic carbon compounds in the C9-C40 carbon range are calculated against the aliphatic calibration curve, and reported as Aliphatic EPH. Therefore, the aliphatic C9-C40 concentration for the sample is reported as the Total EPH.

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Report of Analysis

Client:	PSEG			Date	Collected:	05/07/25		
Project:	OR-636 (Dradell and New Milford		Date	Received:	05/08/25		
Client Sample ID:	OR-636-0)4		SDG	No.:	Q1983		
Lab Sample ID:	Q1983-09)		Matri	x:	Solid		
Analytical Method:	NJEPH			% So	id:	86.8		
Sample Wt/Vol:	30.08	Units: g		Final	Vol:	2000	uL	
Soil Aliquot Vol:		uL		Test:		EPH_NF		
Prep Method :						_		
File ID :	Dilution:	Prep Date :		Date Analy	zed :	P	rep Batch ID	
FC068815.D	1	05/09/25		05/09/25		P	B167926	
CAS Number Para	meter		Conc.	Qualifier	MDL		LOQ / CRQL	Units
TARGETS								
Aliphatic C9-C28		Aliphatic C9-C28	11.0		1.04		4.60	mg/kg
Aliphatic C28-C40		Aliphatic C28-C40	3.31		1.36		2.30	mg/kg
SURROGATES								
3383-33-2		1-chlorooctadecane (SURR)	42.6		40 - 140		85%	SPK: 50
84-15-1		ortho-Terphenyl (SURR)	39.0				78%	SPK: 50



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Quantitation Report For Aliphatic EPH Range.

Lab Sample ID:	Q1983-09	Acq On:	09 May 2025 16:26
Client Sample ID:	OR-636-04	Operator:	YP/AJ
Data file:	FC068815.D	Misc:	
Instrument:	FID_C	ALS Vial:	14
Dilution Factor:	1	Sample Multiplier:	1.00

Compound	R.T.		Response	Conc	highest_standard	Units
Aliphatic C9-C12	3.342	6.642	259172	2.614	300	ug/ml
Aliphatic C12-C16	6.643	10.043	780935	8.583	200	ug/ml
Aliphatic C16-C21	10.044	13.410	9397958	109.584	300	ug/ml
Aliphatic C21-C28	13.411	17.074	1954861	23.527	400	ug/ml
Aliphatic C28-C40	17.075	22.075	3659983	43.251	600	ug/ml
Aliphatic EPH	3.342	22.075	16052909	187.559		ug/ml
ortho-Terphenyl (SURR)	11.711	11.711	4451784	39		ug/ml
1-chlorooctadecane (SURR)	13.144	13.144	3419751	42.63		ug/ml
Aliphatic C9-C28	3.342	17.074	12392926	144.308	1200	ug/ml