

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022982.D
 Acq On : 10 Nov 2024 02:04
 Operator : RC/JU
 Sample : PB164808BL
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK808

Quant Time: Nov 10 03:15:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 08 23:49:48 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	44200	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.340	136	170451	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.204	164	134744	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.010	188	348879	20.000	ng/u1	-0.02
79) Chrysene-d12	21.445	240	416194	20.000	ng/u1	#-0.01
88) Perylene-d12	24.669	264	498897	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6708	7.295	ng/uL	0.00
4) Pyridine-d5	3.552	84	93557	34.245	ng/u1	0.00
7) Phenol-d5	6.793	99	114200	34.947	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	69352	36.122	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.134	132	89814	37.528	ng/u1	0.00
15) 4-Methylphenol-d8	8.299	113	96335	35.824	ng/u1	0.00
21) Nitrobenzene-d5	8.734	128	44611	37.595	ng/u1	0.00
24) 2-Nitrophenol-d4	9.446	143	52750	37.288	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	101762	34.349	ng/u1	0.00
31) 4-Chloroaniline-d4	10.493	131	122317	34.608	ng/u1	0.00
46) Dimethylphthalate-d6	13.610	166	384595	39.602	ng/u1	0.00
49) Acenaphthylene-d8	13.887	160	389681	38.456	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.475	143	49835	34.147	ng/u1	0.00
60) Fluorene-d10	15.222	176	329653	38.717	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.363	200	62997	30.143	ng/u1	0.00
73) Anthracene-d10	17.104	188	579559	39.234	ng/u1	-0.02
81) Pyrene-d10	19.498	212	754276	39.346	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.445	264	981657	41.308	ng/u1	-0.02

Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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