

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
 Data File : BP021842.D
 Acq On : 05 Sep 2024 15:07
 Operator : RC/JU
 Sample : PB163099BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK099

Quant Time: Sep 05 16:06:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 05 05:03:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	285009	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.534	136	1160562	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.404	164	733426	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.222	188	1668842	20.000	ng/u1	0.01
79) Chrysene-d12	21.680	240	1760984	20.000	ng/u1	0.00
88) Perylene-d12	25.086	264	1949186	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	53270	6.205	ng/uL	-0.01
4) Pyridine-d5	3.658	84	707805	28.505	ng/u1	-0.02
7) Phenol-d5	6.910	99	881930	28.995	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.063	67	491266	29.679	ng/u1	-0.03
11) 2-Chlorophenol-d4	7.275	132	579957	29.534	ng/u1	-0.03
15) 4-Methylphenol-d8	8.440	113	641636	27.299	ng/u1	-0.02
21) Nitrobenzene-d5	8.893	128	285399	30.566	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.610	143	289011	30.493	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.151	165	500983	26.823	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.663	131	729311	26.840	ng/u1	-0.02
46) Dimethylphthalate-d6	13.816	166	1726084	30.757	ng/u1	0.00
49) Acenaphthylene-d8	14.098	160	1834289	29.112	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	321686	30.129	ng/u1	0.00
60) Fluorene-d10	15.422	176	1459413	30.932	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.545	200	238073	25.937	ng/u1	0.01
73) Anthracene-d10	17.322	188	2361886	29.821	ng/u1	0.00
81) Pyrene-d10	19.698	212	3010882	29.910	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.839	264	3132594	30.139	ng/u1	-0.02

Target Compounds	Qvalue

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

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