

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102622\
 Data File : BP012326.D
 Acq On : 26 Oct 2022 13:29
 Operator : CG/JU
 Sample : PB148367BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 27 00:22:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 26 23:40:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	217488	20.000	ng/u1	0.00
20) Naphthalene-d8	10.622	136	990951	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.469	164	671642	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1502165	20.000	ng/u1	0.01
79) Chrysene-d12	21.327	240	1687505	20.000	ng/u1	0.01
88) Perylene-d12	23.727	264	1941486	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	29423	5.447	ng/uL	0.00
4) Pyridine-d5	3.675	84	356408	22.970	ng/u1	0.00
7) Phenol-d5	6.987	99	501946	25.732	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	316176	27.151	ng/u1	0.00
11) 2-Chlorophenol-d4	7.346	132	392741	27.570	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	404128	26.194	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	190876	29.703	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	204647	30.682	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	381781	26.412	ng/u1	0.00
31) 4-Chloroaniline-d4	10.769	131	582211	27.051	ng/u1	0.00
46) Dimethylphthalate-d6	13.887	166	1444068	31.408	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	1518344	28.898	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	215968	24.819	ng/u1	0.00
60) Fluorene-d10	15.469	176	1211362	30.773	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	192749	24.645	ng/u1	0.00
73) Anthracene-d10	17.334	188	2066841	31.933	ng/u1	0.01
81) Pyrene-d10	19.569	212	2485811	29.340	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	3005199	31.679	ng/u1	0.02

Target Compounds	Qvalue

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

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