

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011941.D
 Acq On : 05 Oct 2022 21:57
 Operator : CG/JU
 Sample : PB148075BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK075

Quant Time: Oct 05 23:25:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	255950	20.000	ng/u1	0.00
20) Naphthalene-d8	10.687	136	1074660	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	660783	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1395441	20.000	ng/u1	0.00
79) Chrysene-d12	21.368	240	1340393	20.000	ng/u1	0.00
88) Perylene-d12	23.798	264	1353138	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	31712	5.464	ng/uL	0.00
4) Pyridine-d5	3.722	84	393499	23.148	ng/u1	0.00
7) Phenol-d5	7.046	99	557731	25.241	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	314800	25.703	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	468546	26.862	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	435387	23.836	ng/u1	0.00
21) Nitrobenzene-d5	9.046	128	215215	26.292	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	229533	26.209	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	418825	25.856	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	623807	25.260	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	1282660	27.145	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1457829	27.063	ng/u1	0.00
54) 4-Nitrophenol-d4	14.728	143	213900	21.860	ng/u1	0.00
60) Fluorene-d10	15.522	176	1160091	28.270	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	168492	19.949	ng/u1	0.00
73) Anthracene-d10	17.380	188	1893689	29.791	ng/u1	0.00
81) Pyrene-d10	19.616	212	2187925	32.231	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.639	264	2099384	30.993	ng/u1	0.00

Target Compounds Qvalue

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

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