

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100424\
 Data File : BP022241.D
 Acq On : 04 Oct 2024 23:54
 Operator : RC/JU
 Sample : PB163635BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK635

Quant Time: Oct 05 05:15:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 05:25:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.692	152	97831	20.000	ng/u1	0.00
20) Naphthalene-d8	10.445	136	366440	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.304	164	284245	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	680565	20.000	ng/u1	0.00
79) Chrysene-d12	21.527	240	699504	20.000	ng/u1	0.00
88) Perylene-d12	24.803	264	772422	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	13534	5.419	ng/uL	0.00
4) Pyridine-d5	3.628	84	191688	26.850	ng/u1	0.00
7) Phenol-d5	6.875	99	218374	27.288	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.028	67	131354	26.656	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	171512	29.622	ng/u1	0.00
15) 4-Methylphenol-d8	8.387	113	177823	28.042	ng/u1	-0.01
21) Nitrobenzene-d5	8.828	128	81920	29.869	ng/u1	0.00
24) 2-Nitrophenol-d4	9.539	143	96451	31.016	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	188585	27.925	ng/u1	0.00
31) 4-Chloroaniline-d4	10.586	131	228702	28.382	ng/u1	0.00
46) Dimethylphthalate-d6	13.704	166	636970	28.989	ng/u1	0.00
49) Acenaphthylene-d8	13.992	160	682176	29.324	ng/u1	0.00
54) 4-Nitrophenol-d4	14.527	143	75129	19.990	ng/u1	0.01
60) Fluorene-d10	15.310	176	554897	28.550	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.433	200	75351	17.833	ng/u1	0.00
73) Anthracene-d10	17.204	188	925725	29.137	ng/u1	0.00
81) Pyrene-d10	19.580	212	1153407	31.827	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.586	264	1205516	29.673	ng/u1	0.01

Target Compounds Qvalue

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

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