

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022042.D
 Acq On : 26 Sep 2024 00:55
 Operator : RC/JU
 Sample : PB163459BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK459

Quant Time: Sep 26 03:19:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	101646	20.000	ng/u1	0.00
20) Naphthalene-d8	10.463	136	381135	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	282984	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.110	188	694653	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	796113	20.000	ng/u1	0.00
88) Perylene-d12	24.827	264	875161	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	18211	7.018	ng/uL	0.00
4) Pyridine-d5	3.634	84	231610	31.224	ng/u1	0.00
7) Phenol-d5	6.887	99	271899	32.701	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	170804	33.361	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	206301	34.293	ng/u1	0.00
15) 4-Methylphenol-d8	8.405	113	216069	32.794	ng/u1	0.00
21) Nitrobenzene-d5	8.846	128	101879	35.714	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	118660	36.686	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	240302	34.211	ng/u1	0.00
31) 4-Chloroaniline-d4	10.599	131	289173	34.503	ng/u1	0.00
46) Dimethylphthalate-d6	13.722	166	808600	36.965	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	845727	36.516	ng/u1	0.00
54) 4-Nitrophenol-d4	14.516	143	113511	30.337	ng/u1	-0.01
60) Fluorene-d10	15.322	176	700011	36.177	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	125228	29.037	ng/u1	0.00
73) Anthracene-d10	17.210	188	1229709	37.920	ng/u1	0.00
81) Pyrene-d10	19.586	212	1576361	38.219	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.604	264	1701144	36.957	ng/u1	0.00

Target Compounds	Qvalue

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

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