

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090723\
 Data File : BP016991.D
 Acq On : 07 Sep 2023 17:00
 Operator : MA/JU
 Sample : PB155065BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK065

Quant Time: Sep 08 01:11:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	343651	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1460713	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.640	164	906423	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	1943537	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	1531187	20.000	ng/u1	0.00
88) Perylene-d12	24.045	264	1484854	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	70969	7.812	ng/uL	0.00
4) Pyridine-d5	3.758	84	821133	31.631	ng/u1	0.00
7) Phenol-d5	7.122	99	948115	30.199	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	623990	31.774	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	717867	31.593	ng/u1	0.00
15) 4-Methylphenol-d8	8.687	113	735496	30.750	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	355560	31.376	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	388126	31.595	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.416	165	662806	29.225	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	1009192	30.843	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	2241166	33.216	ng/u1	0.00
49) Acenaphthylene-d8	14.340	160	2436446	31.290	ng/u1	0.00
54) 4-Nitrophenol-d4	14.851	143	388250	29.458	ng/u1	0.00
60) Fluorene-d10	15.634	176	1881531	32.570	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	277794	23.751	ng/u1	0.00
73) Anthracene-d10	17.504	188	2987695	33.474	ng/u1	0.00
81) Pyrene-d10	19.739	212	3277075	35.137	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.874	264	2628933	34.899	ng/u1	0.00

Target Compounds Qvalue

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

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