

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018432.D
 Acq On : 28 Nov 2023 10:12
 Operator : MA/JU
 Sample : PB157193BL
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK193

Quant Time: Nov 29 00:59:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 28 08:59:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	434199	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	1868359	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.493	164	1194563	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.245	188	2403970	20.000	ng/u1	0.00
79) Chrysene-d12	21.398	240	1575761	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	1690254	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	68554	5.914	ng/uL	0.00
4) Pyridine-d5	3.681	84	964218	30.266	ng/u1	0.00
7) Phenol-d5	7.022	99	1211747	30.537	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	685242	28.878	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	1012811	30.399	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	1006005	31.415	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	498540	32.852	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	500426	35.706	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	1048084	30.997	ng/u1	0.00
31) 4-Chloroaniline-d4	10.793	131	1521671	33.367	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	3367222	31.679	ng/u1	0.00
49) Acenaphthylene-d8	14.187	160	3724962	30.756	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	381427	25.994	ng/u1	0.00
60) Fluorene-d10	15.487	176	2802853	31.526	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	275483	26.920	ng/u1	0.00
73) Anthracene-d10	17.345	188	4133806	32.172	ng/u1	0.00
81) Pyrene-d10	19.598	212	4156607	33.144	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.804	264	3257197	33.239	ng/u1	0.01

Target Compounds	Qvalue

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

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