

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120723\
 Data File : BP018674.D
 Acq On : 07 Dec 2023 15:45
 Operator : MA/JU
 Sample : PB157560BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK560

Quant Time: Dec 07 22:11:13 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 04 21:39:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	196500	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.640	136	743225	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.475	164	451517	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.222	188	1084831	20.000	ng/u1	0.00
79) Chrysene-d12	21.310	240	1187709	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	1397439	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	36571	6.342	ng/uL	0.00
4) Pyridine-d5	3.681	84	497753	32.227	ng/u1	0.00
7) Phenol-d5	7.016	99	584256	29.628	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	340080	29.098	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.375	132	447052	29.054	ng/u1	-0.01
15) 4-Methylphenol-d8	8.557	113	458075	28.704	ng/u1	-0.01
21) Nitrobenzene-d5	9.004	128	211096	31.912	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.728	143	225377	32.863	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.263	165	433934	30.943	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.781	131	669122	36.510	ng/u1	-0.01
46) Dimethylphthalate-d6	13.892	166	1547710	38.406	ng/u1	-0.01
49) Acenaphthylene-d8	14.169	160	1590637	35.846	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.675	143	285065	43.095	ng/u1	-0.01
60) Fluorene-d10	15.469	176	1294538	38.270	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	218543	33.598	ng/u1	0.00
73) Anthracene-d10	17.322	188	2277875	39.941	ng/u1	0.00
81) Pyrene-d10	19.563	212	2971638	32.142	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	3302286	40.528	ng/u1	0.00

Target Compounds Qvalue

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

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