

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018551.D
 Acq On : 03 Dec 2023 11:51
 Operator : MA/JU
 Sample : PB157194BL
 Misc :
 ALS Vial : 82 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK194

Quant Time: Dec 03 23:42:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 01 21:56:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	364386	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.652	136	1617657	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.487	164	1112556	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	2393117	20.000	ng/u1	0.00
79) Chrysene-d12	21.316	240	1648130	20.000	ng/u1	-0.01
88) Perylene-d12	23.686	264	1667330	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	61925	5.791	ng/uL	0.00
4) Pyridine-d5	3.687	84	897920	31.350	ng/u1	0.00
7) Phenol-d5	7.022	99	1095522	29.959	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.193	67	680010	31.376	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.387	132	832567	29.179	ng/u1	-0.01
15) 4-Methylphenol-d8	8.569	113	893660	30.198	ng/u1	-0.01
21) Nitrobenzene-d5	9.016	128	430069	29.871	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.740	143	450120	30.155	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	879507	28.815	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.793	131	1322471	33.153	ng/u1	-0.01
46) Dimethylphthalate-d6	13.904	166	3130780	31.530	ng/u1	0.00
49) Acenaphthylene-d8	14.181	160	3385772	30.966	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.681	143	459923	28.218	ng/u1	-0.01
60) Fluorene-d10	15.481	176	2599466	31.188	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	342287	23.854	ng/u1	-0.01
73) Anthracene-d10	17.334	188	3994017	31.746	ng/u1	0.00
81) Pyrene-d10	19.569	212	4199352	32.732	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	3088311	31.767	ng/u1	-0.02

Target Compounds Qvalue

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

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