

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP123022\
 Data File : BP013370.D
 Acq On : 30 Dec 2022 13:58
 Operator : CG/JU
 Sample : PB149946BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK946

Quant Time: Dec 30 23:37:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 28 00:16:35 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	352897	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1548406	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.569	164	1101244	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.327	188	2607772	20.000	ng/u1	0.00
79) Chrysene-d12	21.415	240	2449476	20.000	ng/u1	0.00
88) Perylene-d12	23.880	264	2245114	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.316	96	53590	6.493	ng/uL	0.00
4) Pyridine-d5	3.746	84	750209	31.997	ng/u1	0.00
7) Phenol-d5	7.087	99	898633	31.263	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.251	67	604516	34.863	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	719578	31.801	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	731125	31.043	ng/u1	0.00
21) Nitrobenzene-d5	9.098	128	360631	31.700	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	404408	31.575	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	746408	30.003	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	1079029	30.724	ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	2753594	32.535	ng/u1	0.00
49) Acenaphthylene-d8	14.263	160	2885133	31.023	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	364231	23.924	ng/u1	0.00
60) Fluorene-d10	15.569	176	2401199	33.535	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	385057	22.945	ng/u1	0.00
73) Anthracene-d10	17.427	188	3873152	32.921	ng/u1	0.00
81) Pyrene-d10	19.663	212	4700707	33.433	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.721	264	3698638	33.051	ng/u1	-0.01

Target Compounds Qvalue

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

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