

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP123022\
 Data File : BP013367.D
 Acq On : 30 Dec 2022 12:04
 Operator : CG/JU
 Sample : PB149944BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK944

Quant Time: Dec 30 23:35:22 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	389780	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1676093	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.569	164	1166320	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.328	188	2497444	20.000	ng/u1	0.00
79) Chrysene-d12	21.416	240	1716334	20.000	ng/u1	0.00
88) Perylene-d12	23.880	264	1785242	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	56978	6.250	ng/uL	0.00
4) Pyridine-d5	3.746	84	781503	30.178	ng/u1	0.00
7) Phenol-d5	7.087	99	919961	28.977	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	624424	32.604	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	756029	30.250	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	749834	28.824	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	373665	30.343	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	418182	30.163	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	769296	28.568	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	1105308	29.074	ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	2756269	30.749	ng/u1	0.00
49) Acenaphthylene-d8	14.263	160	2941284	29.862	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	318476	19.752	ng/u1	0.00
60) Fluorene-d10	15.563	176	2354588	31.050	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	343506	21.373	ng/u1	0.00
73) Anthracene-d10	17.428	188	3488834	30.965	ng/u1	0.00
81) Pyrene-d10	19.663	212	3593183	36.473	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.715	264	2770709	31.137	ng/u1	-0.02

Target Compounds Qvalue

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

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