

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020623\
 Data File : BP013769.D
 Acq On : 06 Feb 2023 10:33
 Operator : CG/JU
 Sample : PB150652BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK652

Quant Time: Feb 06 23:54:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 03 23:44:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.128	152	238299	20.000	ng/ul	0.00
20) Naphthalene-d8	10.952	136	1024883	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.763	164	770176	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.516	188	1911290	20.000	ng/ul	0.00
79) Chrysene-d12	21.592	240	1978428	20.000	ng/ul	0.00
88) Perylene-d12	24.163	264	1885079	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.476	96	31784	5.655	ng/uL	0.00
4) Pyridine-d5	3.905	84	462527	28.680	ng/ul	0.00
7) Phenol-d5	7.269	99	598021	28.554	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.446	67	356666	29.036	ng/ul	0.00
11) 2-Chlorophenol-d4	7.652	132	436625	28.440	ng/ul	0.00
15) 4-Methylphenol-d8	8.828	113	480713	28.792	ng/ul	0.00
21) Nitrobenzene-d5	9.299	128	198541	30.367	ng/ul	0.00
24) 2-Nitrophenol-d4	10.028	143	238205	31.263	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.563	165	510440	28.648	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.087	131	693527	29.471	ng/ul	0.00
46) Dimethylphthalate-d6	14.163	166	1824622	30.055	ng/ul	0.00
49) Acenaphthylene-d8	14.457	160	1982777	29.560	ng/ul	0.00
54) 4-Nitrophenol-d4	14.946	143	197189	19.270	ng/ul	0.00
60) Fluorene-d10	15.751	176	1658587	31.695	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	164257	15.336	ng/ul	0.00
73) Anthracene-d10	17.616	188	2879864	32.325	ng/ul	0.00
81) Pyrene-d10	19.834	212	3683943	33.386	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.992	264	3142402	32.655	ng/ul	-0.01

Target Compounds Qvalue

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

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