

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013292.D
 Acq On : 24 Dec 2022 10:44
 Operator : CG/JU
 Sample : PB149810BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK810

Quant Time: Dec 25 00:19:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 00:19:17 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	504473	20.000	ng/u1	0.00
20) Naphthalene-d8	10.751	136	1914239	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	1059257	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	2389529	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	2562606	20.000	ng/u1	0.00
88) Perylene-d12	23.921	264	2940441	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	77044	6.530	ng/uL	0.00
4) Pyridine-d5	3.758	84	1046906	31.236	ng/u1	0.00
7) Phenol-d5	7.105	99	1293710	31.484	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	825426	33.300	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	1075166	33.239	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	983110	29.200	ng/u1	0.00
21) Nitrobenzene-d5	9.110	128	496523	35.303	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	553400	34.950	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	968946	31.505	ng/u1	0.00
31) 4-Chloroaniline-d4	10.893	131	1286381	29.628	ng/u1	0.00
46) Dimethylphthalate-d6	13.992	166	2592457	31.845	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	3037598	33.957	ng/u1	0.00
54) 4-Nitrophenol-d4	14.787	143	416245	28.425	ng/u1	0.00
60) Fluorene-d10	15.581	176	2286343	33.197	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	274026	17.820	ng/u1	0.00
73) Anthracene-d10	17.445	188	3824626	35.478	ng/u1	0.00
81) Pyrene-d10	19.680	212	4954138	33.680	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.757	264	5193051	35.432	ng/u1	0.00

Target Compounds Qvalue

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

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