

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012731.D
 Acq On : 25 Nov 2022 17:42
 Operator : CG/JU
 Sample : PB149034BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK034

Quant Time: Nov 26 06:52:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 25 15:59:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	477623	20.000	ng/u1	0.00
20) Naphthalene-d8	10.840	136	2138680	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	1468345	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.416	188	3153139	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	2862673	20.000	ng/u1	0.00
88) Perylene-d12	24.021	264	2505489	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	67765	6.080	ng/uL	0.00
4) Pyridine-d5	3.817	84	940043	28.182	ng/u1	0.00
7) Phenol-d5	7.169	99	1171591	29.851	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	768618	31.782	ng/u1	0.00
11) 2-Chlorophenol-d4	7.552	132	944211	31.573	ng/u1	0.00
15) 4-Methylphenol-d8	8.722	113	969688	30.665	ng/u1	0.00
21) Nitrobenzene-d5	9.193	128	448702	34.278	ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	504417	36.292	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.452	165	991129	30.562	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	1429219	32.183	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	3515831	34.786	ng/u1	0.00
49) Acenaphthylene-d8	14.351	160	3866432	32.787	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	494776	27.458	ng/u1	0.00
60) Fluorene-d10	15.651	176	2994038	33.826	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	397727	26.391	ng/u1	0.00
73) Anthracene-d10	17.516	188	4882082	35.174	ng/u1	0.00
81) Pyrene-d10	19.739	212	5517057	33.168	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.857	264	4326529	35.156	ng/u1	0.00

Target Compounds Qvalue

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

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