

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012762.D
 Acq On : 26 Nov 2022 11:54
 Operator : CG/JU
 Sample : PB149078BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK078

Quant Time: Nov 27 21:50:30 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	484988	20.000	ng/u1	0.00
20) Naphthalene-d8	10.834	136	2101785	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.657	164	1391418	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	2819534	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	1739238	20.000	ng/u1	-0.01
88) Perylene-d12	24.010	264	1452283	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	79638	7.036	ng/uL	0.00
4) Pyridine-d5	3.823	84	1109214	32.749	ng/u1	0.00
7) Phenol-d5	7.164	99	1339164	33.603	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	857699	34.927	ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	1077944	35.497	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	1076980	33.541	ng/u1	-0.01
21) Nitrobenzene-d5	9.187	128	497418	38.667	ng/u1	0.00
24) 2-Nitrophenol-d4	9.911	143	552213	40.428	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	1093035	34.296	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	1553985	35.606	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	3577882	37.358	ng/u1	0.00
49) Acenaphthylene-d8	14.351	160	4039846	36.151	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	450690	26.394	ng/u1	0.00
60) Fluorene-d10	15.646	176	3081180	36.735	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	350824	26.033	ng/u1	0.00
73) Anthracene-d10	17.510	188	4695558	37.833	ng/u1	0.00
81) Pyrene-d10	19.734	212	4748189	46.984	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.845	264	2774143	38.889	ng/u1	-0.01

Target Compounds Qvalue

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

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