

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
 Data File : BP012526.D
 Acq On : 07 Nov 2022 12:04
 Operator : CG/JU
 Sample : PB148749BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK749

Quant Time: Nov 07 21:39:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:53:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	276803	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1325077	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	927110	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	2096666	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	1686717	20.000	ng/ul	0.00
88) Perylene-d12	23.669	264	1482071	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	32828	4.775	ng/uL	0.00
4) Pyridine-d5	3.646	84	167	0.008	ng/ul	0.00
7) Phenol-d5	6.958	99	658602	26.528	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.117	67	373085	25.173	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	497969	27.466	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	543082	27.658	ng/ul	0.00
21) Nitrobenzene-d5	8.952	128	253377	29.487	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	293432	32.900	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	529201	27.379	ng/ul	0.00
31) 4-Chloroaniline-d4	10.705	131	18	0.001	ng/ul	-0.02
46) Dimethylphthalate-d6	13.846	166	1775823	27.981	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	1878287	25.898	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	213780	17.798	ng/ul	0.01
60) Fluorene-d10	15.428	176	1560324	28.715	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	125575	11.504	ng/ul	0.00
73) Anthracene-d10	17.292	188	2469773	27.338	ng/ul	0.00
81) Pyrene-d10	19.534	212	2716532	32.078	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.516	264	2013631	27.806	ng/ul	0.00

Target Compounds Qvalue

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

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