

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012508.D
 Acq On : 04 Nov 2022 02:27
 Operator : CG/JU
 Sample : PB148622BL
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK622

Quant Time: Nov 04 03:47:06 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	318557	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1512544	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.428	164	1070611	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2427159	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	2220589	20.000	ng/u1	0.00
88) Perylene-d12	23.668	264	1967423	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	38727	4.895	ng/uL	0.00
4) Pyridine-d5	3.652	84	276871	12.183	ng/u1	0.00
7) Phenol-d5	6.957	99	762988	26.704	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	433466	25.413	ng/u1	0.00
11) 2-Chlorophenol-d4	7.310	132	575782	27.595	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	634511	28.078	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	298495	30.432	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	346412	34.027	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	625344	28.343	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	539254	16.415	ng/u1	0.00
46) Dimethylphthalate-d6	13.845	166	2239623	30.559	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2408527	28.757	ng/u1	0.00
54) 4-Nitrophenol-d4	14.657	143	358694	25.860	ng/u1	0.00
60) Fluorene-d10	15.428	176	1937053	30.870	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	248898	19.696	ng/u1	0.00
73) Anthracene-d10	17.292	188	3166133	30.274	ng/u1	0.00
81) Pyrene-d10	19.533	212	3600082	32.291	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.515	264	2923742	30.414	ng/u1	0.00

Target Compounds Qvalue

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

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