

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013060.D
 Acq On : 11 Dec 2022 03:15
 Operator : CG/JU
 Sample : PB149498BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 12 03:31:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 01:32:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	646154	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	2845811	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	1974207	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	4452649	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	3941954	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	3813196	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	80742	5.343	ng/uL	0.00
4) Pyridine-d5	3.775	84	1077119	25.090	ng/ul	0.00
7) Phenol-d5	7.122	99	1451907	27.587	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	936578	29.500	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	1192662	28.786	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	1193176	27.668	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	600532	28.721	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	698059	29.655	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1244559	27.220	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	1740938	26.971	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	4433552	29.221	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	4723783	28.333	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	669383	24.526	ng/ul	0.00
60) Fluorene-d10	15.604	176	3740192	29.138	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	715541	24.972	ng/ul	0.00
73) Anthracene-d10	17.463	188	6177726	30.753	ng/ul	0.00
81) Pyrene-d10	19.692	212	7257981	32.077	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.774	264	6097583	32.082	ng/ul	0.00

Target Compounds	Qvalue

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

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