

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012133.D
 Acq On : 14 Oct 2022 12:23
 Operator : CG/JU
 Sample : N4982-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 14 23:02:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	252755	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	1057594	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	621656	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1389280	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1496926	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	1583023	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	25421	3.965	ng/uL	0.00
4) Pyridine-d5	3.705	84	118831	6.970	ng/ul	0.00
7) Phenol-d5	7.016	99	128447	6.319	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	292930	25.387	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	342702	21.238	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	211679	13.557	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	205391	27.498	ng/ul	0.00
24) 2-Nitrophenol-d4	9.734	143	206619	27.281	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	337995	23.027	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	371274	17.165	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1043701	26.510	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	1262748	25.937	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	32808	4.245	ng/ul	0.01
60) Fluorene-d10	15.492	176	990376	27.804	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	143869	19.612	ng/ul	0.00
73) Anthracene-d10	17.351	188	1650949	27.366	ng/ul	0.00
81) Pyrene-d10	19.592	212	2044751	26.663	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.604	264	2116339	27.501	ng/ul	0.00

Target Compounds	Qvalue

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

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