

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012854.D
 Acq On : 29 Nov 2022 16:20
 Operator : CG/JU
 Sample : N5725-17
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 29 21:59:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	469676	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2106589	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1399777	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.404	188	3053238	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	2598489	20.000	ng/ul	-0.01
88) Perylene-d12	23.998	264	2292764	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	39640	3.617	ng/uL	0.00
4) Pyridine-d5	3.823	84	196951	6.004	ng/ul	0.00
7) Phenol-d5	7.158	99	296712	7.688	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	385185	16.197	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	419933	14.279	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	328544	10.566	ng/ul	-0.01
21) Nitrobenzene-d5	9.175	128	203498	15.783	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.905	143	212250	15.503	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	461455	14.446	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	596108	13.627	ng/ul	-0.01
46) Dimethylphthalate-d6	14.051	166	1609384	16.704	ng/ul	0.00
49) Acenaphthylene-d8	14.340	160	1743282	15.507	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	67898	3.953	ng/ul	0.00
60) Fluorene-d10	15.640	176	1468009	17.398	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	137885	9.449	ng/ul	-0.01
73) Anthracene-d10	17.498	188	2427432	18.061	ng/ul	-0.01
81) Pyrene-d10	19.728	212	2794317	18.507	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	1995719	17.721	ng/ul	-0.01

Target Compounds Qvalue

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

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