

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
 Data File : BP012816.D
 Acq On : 28 Nov 2022 14:47
 Operator : CG/JU
 Sample : N5659-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ESQT7

Quant Time: Nov 28 22:39:34 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	8.016	152	337488	20.000	ng/u1	0.00
20) Naphthalene-d8	10.834	136	1486269	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	959810	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	2089600	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	2096576	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	2124692	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	56161	7.131	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.164	99	900951	32.487	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	664921	38.910	ng/u1	0.00
11) 2-Chlorophenol-d4	7.540	132	725940	34.353	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	798675	35.744	ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	354194	38.936	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	354045	36.654	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	736813	32.693	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	955894	30.973	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2735497	41.406	ng/u1	0.00
49) Acenaphthylene-d8	14.346	160	3087535	40.054	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	185793	15.774	ng/u1	0.00
60) Fluorene-d10	15.645	176	2485833	42.964	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	205446	20.571	ng/u1	0.00
73) Anthracene-d10	17.504	188	4001699	43.505	ng/u1	0.00
81) Pyrene-d10	19.733	212	4796447	39.372	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.845	264	4552536	43.622	ng/u1	0.00

Target Compounds Qvalue

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

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