

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
 Data File : BP012817.D
 Acq On : 28 Nov 2022 15:22
 Operator : CG/JU
 Sample : N5659-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ESQT8

Quant Time: Nov 28 22:39:55 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	460870	20.000	ng/u1	0.00
20) Naphthalene-d8	10.834	136	2066849	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1374510	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	3043788	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	3135412	20.000	ng/u1	0.00
88) Perylene-d12	24.010	264	3270315	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	68665	6.384	ng/uL	0.00
4) Pyridine-d5	3.840	84	19075	0.593	ng/u1	0.02
7) Phenol-d5	7.163	99	1315833	34.745	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	853095	36.557	ng/u1	0.00
11) 2-Chlorophenol-d4	7.540	132	1040569	36.059	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	1075221	35.238	ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	491509	38.854	ng/u1	0.00
24) 2-Nitrophenol-d4	9.904	143	542257	40.370	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	1063380	33.929	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1243815	28.981	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	3737006	39.499	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	4156270	37.651	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	527996	31.302	ng/u1	0.00
60) Fluorene-d10	15.645	176	3317580	40.040	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	425605	29.256	ng/u1	0.00
73) Anthracene-d10	17.504	188	5469948	40.825	ng/u1	0.00
81) Pyrene-d10	19.733	212	6830557	37.492	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.845	264	6820542	42.460	ng/u1	0.00

Target Compounds Qvalue

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

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