

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012786.D
 Acq On : 27 Nov 2022 02:07
 Operator : CG/JU
 Sample : N5710-05
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KY4

Quant Time: Nov 27 23:02:21 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	475836	20.000 ng/u1	0.00
20) Naphthalene-d8	10.834	136	2172353	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1486303	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.404	188	3296588	20.000 ng/u1	0.00
79) Chrysene-d12	21.492	240	3233275	20.000 ng/u1	0.00
88) Perylene-d12	24.010	264	3214258	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.393	96	44560	4.013 ng/uL	0.00
4) Pyridine-d5	3.823	84	271598	8.173 ng/u1	0.00
7) Phenol-d5	7.164	99	249750	6.387 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	808835	33.570 ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	765758	25.702 ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	523446	16.615 ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	496005	37.305 ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	481378	34.097 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	1010859	30.687 ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1267582	28.101 ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	4034342	39.434 ng/u1	0.00
49) Acenaphthylene-d8	14.346	160	4220989	35.361 ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	76019	4.168 ng/u1	0.00
60) Fluorene-d10	15.645	176	3459567	38.613 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	277468	17.610 ng/u1	0.00
73) Anthracene-d10	17.504	188	5821019	40.114 ng/u1	0.00
81) Pyrene-d10	19.733	212	6861806	36.524 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.845	264	6413016	40.619 ng/u1	0.00

Target Compounds	Qvalue

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

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