

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012504.D
 Acq On : 03 Nov 2022 23:35
 Operator : CG/JU
 Sample : N5319-15
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAB0

Quant Time: Nov 04 01:04:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	290437	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1371269	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	953931	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2152920	20.000	ng/u1	-0.01
79) Chrysene-d12	21.286	240	2021735	20.000	ng/u1	-0.01
88) Perylene-d12	23.668	264	1781187	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	26968	3.739	ng/uL	0.00
4) Pyridine-d5	3.664	84	120305	5.806	ng/u1	0.00
7) Phenol-d5	6.957	99	142142	5.457	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	471193	30.300	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.310	132	447925	23.546	ng/u1	-0.01
15) 4-Methylphenol-d8	8.499	113	301810	14.649	ng/u1	0.00
21) Nitrobenzene-d5	8.951	128	329351	37.037	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	344257	37.299	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	590147	29.504	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.728	131	755353	25.362	ng/u1	-0.01
46) Dimethylphthalate-d6	13.851	166	2283069	34.962	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2440539	32.704	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.675	143	41284	3.340	ng/u1	0.01
60) Fluorene-d10	15.428	176	1970427	35.243	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.563	200	318075	28.377	ng/u1	-0.01
73) Anthracene-d10	17.292	188	3375514	36.388	ng/u1	-0.01
81) Pyrene-d10	19.533	212	3911010	38.530	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.515	264	3236458	37.187	ng/u1	-0.01

Target Compounds	Qvalue

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

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