

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012350.D
 Acq On : 27 Oct 2022 19:32
 Operator : CG/JU
 Sample : N5015-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C1BF8

Quant Time: Oct 27 23:23:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 26 23:40:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	318544	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.610	136	1432018	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.457	164	971101	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	2090802	20.000	ng/u1	0.00
79) Chrysene-d12	21.315	240	2105592	20.000	ng/u1	0.00
88) Perylene-d12	23.709	264	1820597	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	36315	4.590	ng/uL	0.00
4) Pyridine-d5	3.669	84	493702	21.724	ng/u1	0.00
7) Phenol-d5	6.981	99	731706	25.610	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	447279	26.224	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	566942	27.173	ng/u1	0.00
15) 4-Methylphenol-d8	8.522	113	592590	26.224	ng/u1	-0.01
21) Nitrobenzene-d5	8.975	128	293532	31.609	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.698	143	321529	33.359	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	583141	27.917	ng/u1	0.00
31) 4-Chloroaniline-d4	10.757	131	864935	27.809	ng/u1	0.00
46) Dimethylphthalate-d6	13.875	166	1918128	28.854	ng/u1	0.00
49) Acenaphthylene-d8	14.151	160	2194621	28.888	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	290788	23.112	ng/u1	0.00
60) Fluorene-d10	15.457	176	1696231	29.802	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	223247	20.508	ng/u1	0.00
73) Anthracene-d10	17.321	188	2666436	29.598	ng/u1	0.00
81) Pyrene-d10	19.557	212	3098797	29.313	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.556	264	2613118	29.375	ng/u1	0.00

Target Compounds Qvalue

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

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