

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022365.D
 Acq On : 14 Oct 2024 21:27
 Operator : RC/JU
 Sample : P4237-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4F32

Quant Time: Oct 15 01:35:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 14 11:01:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.698	152	100543	20.000	ng/u1	0.00
20) Naphthalene-d8	10.457	136	386860	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	306804	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	775219	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	811887	20.000	ng/u1	0.00
88) Perylene-d12	24.839	264	980996	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	9563	3.696	ng/uL	0.00
4) Pyridine-d5	3.634	84	141268	19.889	ng/u1	0.00
7) Phenol-d5	6.887	99	202863	23.969	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	116310	23.382	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	154837	25.783	ng/u1	0.00
15) 4-Methylphenol-d8	8.398	113	169888	25.153	ng/u1	0.00
21) Nitrobenzene-d5	8.839	128	76549	25.734	ng/u1	0.00
24) 2-Nitrophenol-d4	9.551	143	93846	27.251	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.086	165	187827	25.662	ng/u1	0.00
31) 4-Chloroaniline-d4	10.592	131	209906	24.270	ng/u1	0.00
46) Dimethylphthalate-d6	13.710	166	650577	26.689	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	708905	27.381	ng/u1	0.00
54) 4-Nitrophenol-d4	14.539	143	90393	22.104	ng/u1	0.00
60) Fluorene-d10	15.316	176	604308	28.582	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	98645	19.348	ng/u1	0.00
73) Anthracene-d10	17.198	188	1016605	27.876	ng/u1	0.00
81) Pyrene-d10	19.586	212	1338538	31.177	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.615	264	1555484	29.690	ng/u1	-0.02

Target Compounds	Qvalue

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

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