

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101124\
 Data File : BP022346.D
 Acq On : 11 Oct 2024 22:58
 Operator : RC/JU
 Sample : P4237-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4EP7

Quant Time: Oct 12 01:08:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 08 02:14:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	96438	20.000	ng/ul	0.00
20) Naphthalene-d8	10.457	136	362817	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	287655	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	731402	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	800181	20.000	ng/ul	0.00
88) Perylene-d12	24.839	264	948112	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	6628	2.671	ng/uL	0.00
4) Pyridine-d5	3.640	84	99168	14.556	ng/ul	0.00
7) Phenol-d5	6.887	99	132500	16.322	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	74478	15.610	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	100492	17.446	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	104741	16.168	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	47823	17.143	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	58236	18.031	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	119651	17.431	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	131114	16.165	ng/ul	0.00
46) Dimethylphthalate-d6	13.716	166	391366	17.124	ng/ul	0.00
49) Acenaphthylene-d8	13.998	160	438149	18.050	ng/ul	0.00
54) 4-Nitrophenol-d4	14.534	143	52183	13.610	ng/ul	0.00
60) Fluorene-d10	15.316	176	369336	18.631	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	51049	10.612	ng/ul	0.00
73) Anthracene-d10	17.210	188	652547	18.966	ng/ul	0.00
81) Pyrene-d10	19.592	212	854097	20.184	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.621	264	987133	19.495	ng/ul	0.02

Target Compounds	Qvalue

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

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