

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022985.D
 Acq On : 10 Nov 2024 04:07
 Operator : RC/JU
 Sample : P4746-16
 Misc :
 ALS Vial : 72 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP89

Quant Time: Nov 10 04:35:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 08 23:49:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	58992	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.334	136	224471	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.216	164	182485	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.010	188	449967	20.000	ng/u1	-0.02
79) Chrysene-d12	21.445	240	550319	20.000	ng/u1	-0.01
88) Perylene-d12	24.668	264	705228	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6893	5.617	ng/uL	0.00
4) Pyridine-d5	3.564	84	18035	4.946	ng/u1	0.00
7) Phenol-d5	6.793	99	76095	17.447	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	88842	34.671	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.134	132	109858	34.393	ng/u1	0.00
15) 4-Methylphenol-d8	8.293	113	109161	30.415	ng/u1	-0.01
21) Nitrobenzene-d5	8.728	128	57306	36.671	ng/u1	0.00
24) 2-Nitrophenol-d4	9.440	143	70312	37.741	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	138575	35.518	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.487	131	59931	12.876	ng/u1	-0.01
46) Dimethylphthalate-d6	13.610	166	504612	38.367	ng/u1	0.00
49) Acenaphthylene-d8	13.898	160	497887	36.280	ng/u1	0.00
54) 4-Nitrophenol-d4	14.498	143	23711	11.996	ng/u1	0.03
60) Fluorene-d10	15.222	176	434080	37.644	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.363	200	91903	34.095	ng/u1	0.00
73) Anthracene-d10	17.110	188	756100	39.686	ng/u1	-0.01
81) Pyrene-d10	19.486	212	990034	39.058	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.451	264	1326450	39.486	ng/u1	-0.02

Target Compounds Qvalue

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

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