

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
 Data File : BP017757.D
 Acq On : 23 Oct 2023 19:27
 Operator : MA/JU
 Sample : 04926-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCCZ6

Quant Time: Oct 24 05:48:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	111218	20.000	ng/u1	0.00
20) Naphthalene-d8	10.716	136	444197	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	267178	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.375	188	558009	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.516	240	444000	20.000	ng/u1	0.00
88) Perylene-d12	24.069	264	456079	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	15397	5.150	ng/uL	0.00
4) Pyridine-d5	3.687	84	139404	17.810	ng/u1	0.00
7) Phenol-d5	7.046	99	255318	27.587	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	177254	31.012	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	215115	28.498	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	166469	22.905	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	94421	26.233	ng/u1	0.00
24) 2-Nitrophenol-d4	9.810	143	97984	24.638	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	194704	24.958	ng/u1	0.00
31) 4-Chloroaniline-d4	10.893	131	188778	18.491	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	650210	28.990	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	734854	29.379	ng/u1	0.00
54) 4-Nitrophenol-d4	14.840	143	47093	14.736	ng/u1	0.01
60) Fluorene-d10	15.587	176	557427	28.751	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	55742	15.166	ng/u1	0.00
73) Anthracene-d10	17.475	188	818923	28.718	ng/u1	0.00
81) Pyrene-d10	19.733	212	957572	34.400	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	758231	29.849	ng/u1	0.00

Target Compounds	Qvalue

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

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