

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
 Data File : BP017756.D
 Acq On : 23 Oct 2023 18:53
 Operator : MA/JU
 Sample : 04926-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCCZ2

Quant Time: Oct 24 05:42:28 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	121996	20.000	ng/u1	0.00
20) Naphthalene-d8	10.716	136	485935	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	291686	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.375	188	598620	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.516	240	468870	20.000	ng/u1	0.00
88) Perylene-d12	24.068	264	489559	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	13544	4.130	ng/uL	0.00
4) Pyridine-d5	3.693	84	150951	17.581	ng/u1	0.00
7) Phenol-d5	7.046	99	246695	24.300	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	168014	26.798	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	204949	24.752	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	179041	22.458	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	87863	22.314	ng/u1	0.00
24) 2-Nitrophenol-d4	9.810	143	82588	18.983	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	179827	21.071	ng/u1	0.00
31) 4-Chloroaniline-d4	10.899	131	122699	10.986	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	592710	24.206	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	674303	24.693	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	41991	12.035	ng/u1	0.01
60) Fluorene-d10	15.587	176	515085	24.335	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	47196	11.969	ng/u1	0.00
73) Anthracene-d10	17.475	188	749913	24.514	ng/u1	0.00
81) Pyrene-d10	19.733	212	863668	29.381	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	688024	25.233	ng/u1	0.00

Target Compounds Qvalue

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

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