

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\
 Data File : BG049242.D
 Acq On : 9 Jul 2021 18:00
 Operator : CG/JU
 Sample : PB137522BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jul 09 18:57:17 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 09 03:15:38 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	27104	20.00	ng/ul	0.00
20) Naphthalene-d8	10.89	136	112612	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.72	164	79725	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.47	188	192572	20.00	ng/ul	0.00
79) Chrysene-d12	21.75	240	200259	20.00	ng/ul	0.00
88) Perylene-d12	25.05	264	200169	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	4645m	8.06	ng/uL	0.00
4) Pyridine-d5	3.89	84	55674	32.86	ng/ul	0.00
7) Phenol-d5	7.23	99	80600	34.19	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.40	67	50998	37.29	ng/ul	0.00
11) 2-Chlorophenol-d4	7.60	132	62161	35.67	ng/ul	0.00
15) 4-Methylphenol-d8	8.78	113	66541	35.23	ng/ul	0.00
21) Nitrobenzene-d5	9.24	128	32473	37.58	ng/ul	0.00
24) 2-Nitrophenol-d4	9.97	143	37710	38.01	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.51	165	69502	35.91	ng/ul	0.00
31) 4-Chloroaniline-d4	11.02	131	89025	35.29	ng/ul	0.00
46) Dimethylphthalate-d6	14.12	166	246119	40.14	ng/ul	0.00
49) Acenaphthylene-d8	14.41	160	278267	38.68	ng/ul	0.00
54) 4-Nitrophenol-d4	14.91	143	36116	35.59	ng/ul	0.00
60) Fluorene-d10	15.71	176	201531	38.82	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.82	200	41909	34.24	ng/ul	0.00
73) Anthracene-d10	17.57	188	339401	38.71	ng/ul	0.00
81) Pyrene-d10	19.85	212	416828	40.61	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.82	264	412552	39.65	ng/ul	0.00

Target Compounds Ovalue

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

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