

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062022\
 Data File : BG053959.D
 Acq On : 20 Jun 2022 21:54
 Operator : CG/JU
 Sample : N3358-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW4S2

Quant Time: Jun 20 23:59:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 20 15:29:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.080	152	19786	20.000	ng/ul	0.00
20) Naphthalene-d8	10.888	136	79368	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.702	164	60511	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	162685	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.717	240	183700	20.000	ng/ul	# 0.00
88) Perylene-d12	24.978	264	187487	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.503	96	2149	5.168	ng/uL	0.00
4) Pyridine-d5	3.920	84	10254	9.066	ng/ul	0.00
7) Phenol-d5	7.228	99	12345	8.709	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.398	67	29249	34.207	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	32060	28.363	ng/ul	0.00
15) 4-Methylphenol-d8	8.773	113	23336	19.308	ng/ul	0.00
21) Nitrobenzene-d5	9.237	128	20346	34.537	ng/ul	0.00
24) 2-Nitrophenol-d4	9.966	143	22114	35.329	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.501	165	46547	33.611	ng/ul	0.00
31) 4-Chloroaniline-d4	11.018	131	24951	14.261	ng/ul	0.00
46) Dimethylphthalate-d6	14.102	166	186939	39.270	ng/ul	0.00
49) Acenaphthylene-d8	14.396	160	192066	37.105	ng/ul	0.00
54) 4-Nitrophenol-d4	14.878	143	5183	7.952	ng/ul	0.00
60) Fluorene-d10	15.689	176	164207	38.403	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.800	200	33488	37.735	ng/ul	0.00
73) Anthracene-d10	17.545	188	285475	38.855	ng/ul	0.00
81) Pyrene-d10	19.825	212	385996	40.300	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.754	264	373513	39.467	ng/ul	0.00

Target Compounds Qvalue

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

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