

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\
 Data File : BM033139.D
 Acq On : 18 Nov 2021 02:59
 Operator : CG/JU
 Sample : M4618-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG218

Quant Time: Nov 18 03:35:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 17 14:14:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.972	152	83810	20.000	ng/u1	0.00
20) Naphthalene-d8	10.772	136	328613	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.589	164	211117	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.324	188	438838	20.000	ng/u1	0.00
79) Chrysene-d12	21.477	240	434171	20.000	ng/u1	0.00
88) Perylene-d12	23.824	264	445798	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.425	96	9647	4.438	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.125	99	38589	5.443	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.301	67	98184	21.835	ng/u1	0.00
11) 2-Chlorophenol-d4	7.501	132	94755	17.647	ng/u1	0.00
15) 4-Methylphenol-d8	8.666	113	66014	11.999	ng/u1	0.00
21) Nitrobenzene-d5	9.131	128	54556	23.124	ng/u1	0.00
24) 2-Nitrophenol-d4	9.848	143	52852	22.368	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.383	165	109401	20.218	ng/u1	0.00
31) 4-Chloroaniline-d4	10.901	131	89145	12.388	ng/u1	0.00
46) Dimethylphthalate-d6	13.995	166	398699	25.747	ng/u1	-0.01
49) Acenaphthylene-d8	14.283	160	463482	23.233	ng/u1	0.00
54) 4-Nitrophenol-d4	14.765	143	13098	5.193	ng/u1	0.00
60) Fluorene-d10	15.577	176	351275	25.288	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.683	200	46849	22.716	ng/u1	-0.01
73) Anthracene-d10	17.424	188	638769	30.196	ng/u1	0.00
81) Pyrene-d10	19.700	212	744295	29.009	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.671	264	714737	29.874	ng/u1	-0.01

Target Compounds Qvalue

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

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