

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
 Data File : BM042905.D
 Acq On : 19 Nov 2023 14:24
 Operator : MA/JU
 Sample : 05370-16
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCDR0

Quant Time: Nov 19 21:49:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 18 22:06:09 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	28559	20.000	ng/u1	0.00
20) Naphthalene-d8	10.622	136	113457	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.468	164	72952	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	164008	20.000	ng/u1	0.00
79) Chrysene-d12	21.409	240	169887	20.000	ng/u1	0.00
88) Perylene-d12	23.762	264	206469	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	3884	4.267	ng/uL	0.00
4) Pyridine-d5	3.616	84	19639	8.675	ng/u1	0.01
7) Phenol-d5	6.987	99	11350	4.083	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.163	67	59209	30.315	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	41184	19.203	ng/u1	0.00
15) 4-Methylphenol-d8	8.539	113	22647	10.348	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	23044	23.166	ng/u1	0.00
24) 2-Nitrophenol-d4	9.728	143	26465	22.743	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	46250	21.375	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	48384	17.612	ng/u1	0.00
46) Dimethylphthalate-d6	13.892	166	190667	27.615	ng/u1	0.00
49) Acenaphthylene-d8	14.168	160	194927	26.386	ng/u1	0.00
54) 4-Nitrophenol-d4	14.792	143	77	0.099	ng/u1	0.07
60) Fluorene-d10	15.462	176	157032	27.466	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.615	200	21930	17.653	ng/u1	0.00
73) Anthracene-d10	17.321	188	271274	29.514	ng/u1	0.00
81) Pyrene-d10	19.621	212	372083	33.446	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.609	264	392955	31.818	ng/u1	0.00

Target Compounds	Qvalue

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

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