

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032435.D
 Acq On : 12 Oct 2021 00:15
 Operator : CG/JU
 Sample : M4029-17
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00S8

Quant Time: Oct 12 01:03:32 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 11 11:06:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.613	152	216280	20.000	ng/ul	0.00
20) Naphthalene-d8	10.401	136	972750	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.260	164	690549	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.989	188	1446692	20.000	ng/ul	0.00
79) Chrysene-d12	21.153	240	1449713	20.000	ng/ul	0.00
88) Perylene-d12	23.294	264	1527113	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.954	96	9840	1.824	ng/uL	0.00
4) Pyridine-d5	3.396	84	67818	4.450	ng/ul	0.01
7) Phenol-d5	6.801	99	180263	9.802	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.937	67	106758	10.110	ng/ul	0.00
11) 2-Chlorophenol-d4	7.148	132	144215	10.270	ng/ul	0.00
15) 4-Methylphenol-d8	8.342	113	141704	9.539	ng/ul	0.00
21) Nitrobenzene-d5	8.766	128	67695	9.863	ng/ul	0.00
24) 2-Nitrophenol-d4	9.489	143	73460	9.930	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.031	165	141478	9.607	ng/ul	0.00
31) 4-Chloroaniline-d4	10.536	131	190609	8.775	ng/ul	0.00
46) Dimethylphthalate-d6	13.666	166	534241	11.179	ng/ul	0.00
49) Acenaphthylene-d8	13.948	160	588072	10.129	ng/ul	0.00
54) 4-Nitrophenol-d4	14.466	143	67422	7.184	ng/ul	-0.01
60) Fluorene-d10	15.248	176	452079	11.154	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.365	200	50549	6.131	ng/ul	0.00
73) Anthracene-d10	17.083	188	726102	11.530	ng/ul	0.00
81) Pyrene-d10	19.371	212	910838	12.489	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.159	264	899642	11.764	ng/ul	-0.01

Target Compounds	Qvalue

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

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