

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032429.D
 Acq On : 11 Oct 2021 20:38
 Operator : CG/JU
 Sample : M4029-15
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00S6

Quant Time: Oct 12 01:02:08 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	192302	20.000	ng/ul	0.00
20) Naphthalene-d8	10.401	136	877883	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.260	164	630259	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.989	188	1348593	20.000	ng/ul	0.00
79) Chrysene-d12	21.154	240	1379410	20.000	ng/ul	0.00
88) Perylene-d12	23.301	264	1479900	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	10766	2.245	ng/uL	0.00
4) Pyridine-d5	3.396	84	64838	4.785	ng/ul	0.01
7) Phenol-d5	6.802	99	177239	10.839	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.937	67	108172	11.521	ng/ul	0.00
11) 2-Chlorophenol-d4	7.155	132	144872	11.603	ng/ul	0.00
15) 4-Methylphenol-d8	8.343	113	147112	11.138	ng/ul	0.00
21) Nitrobenzene-d5	8.766	128	71703	11.576	ng/ul	0.00
24) 2-Nitrophenol-d4	9.490	143	80402	12.042	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.037	165	143708	10.813	ng/ul	0.00
31) 4-Chloroaniline-d4	10.537	131	211011	10.764	ng/ul	0.00
46) Dimethylphthalate-d6	13.666	166	560147	12.842	ng/ul	0.00
49) Acenaphthylene-d8	13.948	160	635721	11.997	ng/ul	0.00
54) 4-Nitrophenol-d4	14.466	143	62745	7.325	ng/ul	-0.01
60) Fluorene-d10	15.248	176	478116	12.925	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.366	200	49424	6.431	ng/ul	0.00
73) Anthracene-d10	17.083	188	755645	12.872	ng/ul	0.00
81) Pyrene-d10	19.371	212	927028	13.359	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.165	264	951091	12.833	ng/ul	0.00

Target Compounds Qvalue

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

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