

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
 Data File : BP006651.D
 Acq On : 11 Aug 2021 12:11
 Operator : CG/JU
 Sample : M3246-01 30X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA6

Quant Time: Aug 11 11:52:01 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 10 15:09:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	143877	20.000	ng/ul	0.00
20) Naphthalene-d8	10.498	136	619621	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.357	164	353662	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	694242	20.000	ng/ul	0.00
79) Chrysene-d12	21.215	240	651381	20.000	ng/ul	0.00
88) Perylene-d12	23.527	264	675340	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	783	0.184	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.893	99	10088	0.697	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	7707	0.867	ng/ul	0.00
11) 2-Chlorophenol-d4	7.246	132	8741	0.811	ng/ul	-0.01
15) 4-Methylphenol-d8	8.428	113	6617	0.573	ng/ul	-0.01
21) Nitrobenzene-d5	8.881	128	4563	0.868	ng/ul	0.00
24) 2-Nitrophenol-d4	9.593	143	3929	0.714	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	6512	0.700	ng/ul	0.00
31) 4-Chloroaniline-d4	10.651	131	2808	0.186	ng/ul	0.00
46) Dimethylphthalate-d6	13.781	166	22995	0.848	ng/ul	0.00
49) Acenaphthylene-d8	14.051	160	26768	0.805	ng/ul	0.00
54) 4-Nitrophenol-d4	14.575	143	1649	0.291	ng/ul	0.00
60) Fluorene-d10	15.351	176	19101	0.866	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	1338	0.292	ng/ul	0.00
73) Anthracene-d10	17.210	188	29692	0.904	ng/ul	0.00
81) Pyrene-d10	19.457	212	33106	0.968	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.380	264	32339	0.888	ng/ul	0.00

Target Compounds Qvalue

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

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