

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
 Data File : BP006653.D
 Acq On : 11 Aug 2021 13:19
 Operator : CG/JU
 Sample : M3246-06 30X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AB0

Manual Integrations
 APPROVED

mohammad
 8/11/2021 9:29:42 PM

Quant Time: Aug 11 13:17:14 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	145563	20.000	ng/u1	0.00
20) Naphthalene-d8	10.499	136	629084	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	359392	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.110	188	693843	20.000	ng/u1	0.00
79) Chrysene-d12	21.216	240	652188	20.000	ng/u1	0.00
88) Perylene-d12	23.527	264	674505	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	998m	0.232	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	6.905	99	6301	0.431	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	5294	0.589	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	5637	0.517	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	3981	0.341	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	3046	0.570	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	2551	0.457	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	4659	0.494	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	4940	0.322	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	15818	0.574	ng/u1	0.00
49) Acenaphthylene-d8	14.045	160	18433	0.545	ng/u1	0.00
54) 4-Nitrophenol-d4	14.587	143	650	0.113	ng/u1	0.01
60) Fluorene-d10	15.351	176	13630	0.608	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	851m	0.186	ng/u1	-0.01
73) Anthracene-d10	17.210	188	21329	0.650	ng/u1	0.00
81) Pyrene-d10	19.457	212	23606	0.689	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.374	264	23186	0.637	ng/u1	0.00

Target Compounds Qvalue

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

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