

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081021\
 Data File : BP006649.D
 Acq On : 11 Aug 2021 10:21
 Operator : CG/JU
 Sample : PB138246BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK246

Quant Time: Aug 11 11:42:36 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	7.711	152	147595	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.499	136	632127	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	350158	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.116	188	688284	20.000	ng/u1	0.00
79) Chrysene-d12	21.222	240	676422	20.000	ng/u1	0.00
88) Perylene-d12	23.545	264	702369	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	31387	7.193	ng/uL	0.00
4) Pyridine-d5	3.628	84	440021	35.078	ng/u1	0.00
7) Phenol-d5	6.893	99	499523	33.666	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	320692	35.169	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.246	132	389912	35.284	ng/u1	-0.01
15) 4-Methylphenol-d8	8.428	113	399455	33.699	ng/u1	-0.01
21) Nitrobenzene-d5	8.869	128	197868	36.875	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.587	143	207110	36.887	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.122	165	320718	33.809	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.646	131	522407	33.882	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	989661	36.854	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	1250363	37.961	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	182903	32.608	ng/u1	0.00
60) Fluorene-d10	15.357	176	812113	37.185	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.487	200	122802	26.992	ng/u1	0.00
73) Anthracene-d10	17.216	188	1299378	39.900	ng/u1	0.00
81) Pyrene-d10	19.463	212	1418421	39.931	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.392	264	1429841	37.751	ng/u1	0.01

Target Compounds Qvalue

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

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