

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
 Data File : BP006578.D
 Acq On : 07 Aug 2021 21:13
 Operator : CG/JU
 Sample : PB138110BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK110

Quant Time: Aug 08 02:37:05 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Aug 07 13:59:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.752	152	190725	20.000	ng/u1	0.00
20) Naphthalene-d8	10.528	136	827734	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	473208	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.128	188	937262	20.000	ng/u1	0.00
79) Chrysene-d12	21.228	240	873738	20.000	ng/u1	0.00
88) Perylene-d12	23.551	264	913330	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	41456	8.114	ng/uL	0.00
4) Pyridine-d5	3.676	84	589816	38.885	ng/u1	0.00
7) Phenol-d5	6.928	99	654009	36.511	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.093	67	436712	39.183	ng/u1	0.00
11) 2-Chlorophenol-d4	7.281	132	510108	38.331	ng/u1	0.00
15) 4-Methylphenol-d8	8.458	113	520892	36.978	ng/u1	0.00
21) Nitrobenzene-d5	8.905	128	256681	41.647	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	261334	44.408	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.152	165	413634	35.869	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	696039	36.662	ng/u1	0.00
46) Dimethylphthalate-d6	13.799	166	1416773	42.067	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	1722765	41.635	ng/u1	0.00
54) 4-Nitrophenol-d4	14.587	143	237113	35.457	ng/u1	0.00
60) Fluorene-d10	15.375	176	1142280	40.345	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.493	200	170404	32.613	ng/u1	0.00
73) Anthracene-d10	17.228	188	1793621	42.393	ng/u1	0.00
81) Pyrene-d10	19.469	212	1928035	42.902	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	1934657	41.684	ng/u1	0.00

Target Compounds Qvalue

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

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