

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
 Data File : BP006592.D
 Acq On : 08 Aug 2021 06:19
 Operator : CG/JU
 Sample : PB138110BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK110

Quant Time: Aug 09 01:37:35 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	244531	20.000	ng/u1	0.00
20) Naphthalene-d8	10.528	136	1053787	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	585150	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.128	188	1110776	20.000	ng/u1	0.00
79) Chrysene-d12	21.228	240	1036713	20.000	ng/u1	0.00
88) Perylene-d12	23.551	264	1071560	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	49645	7.578	ng/uL	0.00
4) Pyridine-d5	3.681	84	730792	37.578	ng/u1	0.00
7) Phenol-d5	6.928	99	850496	37.033	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.093	67	557126	38.988	ng/u1	0.00
11) 2-Chlorophenol-d4	7.281	132	667898	39.145	ng/u1	0.00
15) 4-Methylphenol-d8	8.458	113	672548	37.239	ng/u1	0.00
21) Nitrobenzene-d5	8.905	128	338575	43.150	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	346561	46.257	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.152	165	543241	37.002	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	891428	36.881	ng/u1	0.00
46) Dimethylphthalate-d6	13.799	166	1728156	41.496	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2171983	42.449	ng/u1	0.00
54) 4-Nitrophenol-d4	14.587	143	307031	37.130	ng/u1	0.00
60) Fluorene-d10	15.375	176	1397239	39.909	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.498	200	207611	33.527	ng/u1	0.00
73) Anthracene-d10	17.228	188	2158128	43.040	ng/u1	0.00
81) Pyrene-d10	19.475	212	2269901	42.569	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	2285002	41.962	ng/u1	0.00

Target Compounds Qvalue

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

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