

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007670.D
 Acq On : 26 Oct 2021 08:23
 Operator : CG/JU
 Sample : PB140163BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK163

Quant Time: Oct 26 09:25:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 25 16:57:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	233773	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	932254	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	606362	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.328	188	1314018	20.000	ng/u1	0.00
79) Chrysene-d12	21.416	240	1465274	20.000	ng/u1	0.00
88) Perylene-d12	23.863	264	1573061	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	34526	6.709	ng/uL	0.00
4) Pyridine-d5	3.781	84	487108	31.300	ng/u1	0.00
7) Phenol-d5	7.099	99	584378	32.911	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	339367	34.424	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	473965	34.320	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	463381	33.328	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	223724	35.215	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	245887	35.437	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	453961	32.142	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.875	131	632733	33.881	ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	1468241	35.747	ng/u1	-0.01
49) Acenaphthylene-d8	14.263	160	1725970	34.285	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	250230	33.236	ng/u1	-0.01
60) Fluorene-d10	15.569	176	1218233	35.182	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	195738	27.848	ng/u1	0.00
73) Anthracene-d10	17.428	188	1989914	36.850	ng/u1	0.00
81) Pyrene-d10	19.663	212	2437679	35.733	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.710	264	2604825	33.708	ng/u1	0.00

Target Compounds Qvalue

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

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