

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007653.D
 Acq On : 25 Oct 2021 22:13
 Operator : CG/JU
 Sample : PB140162BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK162

Quant Time: Oct 26 02:38:52 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	7.940	152	167820	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	683562	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	449930	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.328	188	1001821	20.000	ng/u1	0.00
79) Chrysene-d12	21.416	240	1105013	20.000	ng/u1	0.00
88) Perylene-d12	23.863	264	1172079	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	27821	7.531	ng/uL	0.00
4) Pyridine-d5	3.787	84	423186	37.879	ng/u1	0.00
7) Phenol-d5	7.105	99	481833	37.800	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	280717	39.666	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	383709	38.704	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	377663	37.837	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	181882	39.045	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	194400	38.209	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	370117	35.739	ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	521464	38.082	ng/u1	0.00
46) Dimethylphthalate-d6	13.987	166	1239516	40.670	ng/u1	0.00
49) Acenaphthylene-d8	14.269	160	1420815	38.036	ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	204006	36.517	ng/u1	0.00
60) Fluorene-d10	15.569	176	1037930	40.397	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	156358	29.178	ng/u1	0.00
73) Anthracene-d10	17.428	188	1675465	40.695	ng/u1	0.00
81) Pyrene-d10	19.663	212	2035399	39.564	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.704	264	2172692	37.735	ng/u1	0.00

Target Compounds	Qvalue

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

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