

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007534.D
 Acq On : 20 Oct 2021 18:07
 Operator : CG/JU
 Sample : PB139973BL
 Misc :
 ALS Vial : 89 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK973

Quant Time: Oct 21 01:01:49 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 20 10:49:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	204177	20.00	ng/ul	0.00
20) Naphthalene-d8	10.758	136	839107	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	542252	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.340	188	1163589	20.00	ng/ul	0.00
79) Chrysene-d12	21.422	240	1212184	20.00	ng/ul	0.00
88) Perylene-d12	23.869	264	1247830	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	28717	5.55	ng/uL	0.00
4) Pyridine-d5	3.793	84	415637	29.37	ng/ul	0.00
7) Phenol-d5	7.111	99	494559	31.02	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	283487	29.81	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	394704	31.47	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	391373	31.87	ng/ul	0.00
21) Nitrobenzene-d5	9.111	128	185854	35.00	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	192402	38.79	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	382456	31.61	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	523815	32.02	ng/ul	0.00
46) Dimethylphthalate-d6	13.999	166	1166070	31.08	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1429292	30.32	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	192568	31.99	ng/ul	0.00
60) Fluorene-d10	15.581	176	1002352	31.93	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.693	200	123839	26.56	ng/ul	0.00
73) Anthracene-d10	17.440	188	1616724	32.72	ng/ul	0.00
81) Pyrene-d10	19.669	212	1922530	31.73	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.716	264	1916385	30.86	ng/ul	0.00

Target Compounds	Qvalue

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

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