

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081221\
 Data File : BP006663.D
 Acq On : 12 Aug 2021 13:52
 Operator : CG/JU
 Sample : PB138304BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK304

Quant Time: Aug 12 13:43:13 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 12 13:01:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	139478	20.000	ng/ul	0.00
20) Naphthalene-d8	10.493	136	584432	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.351	164	345197	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	709551	20.000	ng/ul	0.00
79) Chrysene-d12	21.216	240	685641	20.000	ng/ul	0.00
88) Perylene-d12	23.527	264	732465	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	29047	7.044	ng/uL	0.00
4) Pyridine-d5	3.628	84	408420	34.453	ng/ul	0.00
7) Phenol-d5	6.893	99	451548	32.204	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	285176	33.094	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	358530	34.332	ng/ul	0.00
15) 4-Methylphenol-d8	8.422	113	357683	31.931	ng/ul	0.00
21) Nitrobenzene-d5	8.869	128	178801	36.041	ng/ul	0.00
24) 2-Nitrophenol-d4	9.587	143	189423	36.490	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	290542	33.127	ng/ul	0.00
31) 4-Chloroaniline-d4	10.640	131	479530	33.639	ng/ul	0.00
46) Dimethylphthalate-d6	13.775	166	1030594	38.930	ng/ul	0.00
49) Acenaphthylene-d8	14.040	160	1194380	36.782	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143	185509	33.548	ng/ul	0.00
60) Fluorene-d10	15.351	176	804231	37.353	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	135292	28.846	ng/ul	0.00
73) Anthracene-d10	17.210	188	1313793	39.133	ng/ul	0.00
81) Pyrene-d10	19.457	212	1456259	40.445	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.380	264	1491606	37.764	ng/ul	0.00

Target Compounds	Qvalue

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

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