

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030122\
 Data File : BM034101.D
 Acq On : 01 Mar 2022 10:39
 Operator : CG/JU
 Sample : PB142783BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK783

Quant Time: Mar 01 11:51:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 01 03:48:18 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.995	152	136090	20.000	ng/u1	0.00
20) Naphthalene-d8	10.813	136	492243	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.630	164	321094	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.371	188	679997	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.542	240	709973	20.000	ng/u1	# 0.00
88) Perylene-d12	23.994	264	755671	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	18535	6.280	ng/uL	0.00
4) Pyridine-d5	3.790	84	246837	31.826	ng/u1	0.00
7) Phenol-d5	7.142	99	282700	29.086	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.313	67	163416	30.285	ng/u1	0.00
11) 2-Chlorophenol-d4	7.525	132	256903	30.660	ng/u1	0.00
15) 4-Methylphenol-d8	8.695	113	233939	28.278	ng/u1	0.00
21) Nitrobenzene-d5	9.154	128	117458	31.590	ng/u1	0.00
24) 2-Nitrophenol-d4	9.883	143	134741	31.881	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.425	165	254503	29.756	ng/u1	0.00
31) 4-Chloroaniline-d4	10.936	131	320210	29.687	ng/u1	0.00
46) Dimethylphthalate-d6	14.036	166	812410	33.440	ng/u1	0.00
49) Acenaphthylene-d8	14.324	160	950905	32.137	ng/u1	0.00
54) 4-Nitrophenol-d4	14.807	143	115164	27.675	ng/u1	0.00
60) Fluorene-d10	15.618	176	674410	32.117	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.724	200	113402	25.544	ng/u1	0.00
73) Anthracene-d10	17.471	188	1078147	32.788	ng/u1	0.00
81) Pyrene-d10	19.759	212	1242331	32.460	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.836	264	1280178	33.242	ng/u1	0.00

Target Compounds Qvalue

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

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