

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
 Data File : BM034003.D
 Acq On : 03 Feb 2022 15:48
 Operator : CG/JU
 Sample : PB142441BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK441

Quant Time: Feb 04 00:07:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 01 02:16:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.919	152	108095	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.731	136	463725	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.560	164	298215	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.295	188	583346	20.000	ng/u1	-0.01
79) Chrysene-d12	21.459	240	469187	20.000	ng/u1	0.00
88) Perylene-d12	23.847	264	432578	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	15943	5.817	ng/uL	0.00
4) Pyridine-d5	3.713	84	213735	27.939	ng/u1	0.00
7) Phenol-d5	7.078	99	272279	30.067	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.243	67	177582	32.165	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.449	132	228292	31.465	ng/u1	-0.01
15) 4-Methylphenol-d8	8.631	113	227585	31.715	ng/u1	-0.01
21) Nitrobenzene-d5	9.084	128	114774	30.846	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.813	143	126350	31.445	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.348	165	213914	29.419	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.866	131	314033	31.274	ng/u1	0.00
46) Dimethylphthalate-d6	13.966	166	756837	33.962	ng/u1	-0.01
49) Acenaphthylene-d8	14.254	160	872385	31.957	ng/u1	0.00
54) 4-Nitrophenol-d4	14.748	143	89006	23.378	ng/u1	0.00
60) Fluorene-d10	15.548	176	604622	32.493	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.660	200	87786	23.971	ng/u1	-0.01
73) Anthracene-d10	17.395	188	910293	32.584	ng/u1	0.00
81) Pyrene-d10	19.677	212	947479	34.205	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.695	264	752251	33.109	ng/u1	-0.01

Target Compounds Qvalue

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

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