

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092521\
 Data File : BM032181.D
 Acq On : 26 Sep 2021 08:40
 Operator : CG/JU
 Sample : PB139378BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK378

Quant Time: Sep 27 04:10:45 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM092421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 25 01:15:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.660	152	362940	20.000	ng/u1	0.00
20) Naphthalene-d8	10.454	136	1409460	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.307	164	799138	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.036	188	1519609	20.000	ng/u1	0.00
79) Chrysene-d12	21.194	240	1362727	20.000	ng/u1	0.00
88) Perylene-d12	23.359	264	1388019	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.990	96	59677	6.943	ng/uL	0.00
4) Pyridine-d5	3.413	84	807300	33.084	ng/u1	0.00
7) Phenol-d5	6.842	99	962090	33.395	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.984	67	581179	33.550	ng/u1	0.00
11) 2-Chlorophenol-d4	7.201	132	782802	34.999	ng/u1	0.00
15) 4-Methylphenol-d8	8.383	113	740452	32.869	ng/u1	-0.01
21) Nitrobenzene-d5	8.813	128	339756	40.359	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.542	143	355020	39.883	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.083	165	697494	34.276	ng/u1	0.00
31) 4-Chloroaniline-d4	10.589	131	993560	36.125	ng/u1	0.00
46) Dimethylphthalate-d6	13.713	166	1847228	35.577	ng/u1	0.00
49) Acenaphthylene-d8	14.001	160	2372557	35.966	ng/u1	0.00
54) 4-Nitrophenol-d4	14.507	143	261037	30.905	ng/u1	0.00
60) Fluorene-d10	15.295	176	1627250	36.729	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.407	200	181163	25.226	ng/u1	-0.01
73) Anthracene-d10	17.130	188	2373190	37.428	ng/u1	0.00
81) Pyrene-d10	19.418	212	2556159	35.651	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.224	264	2323772	35.673	ng/u1	-0.01

Target Compounds	Qvalue

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

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