

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092521\
 Data File : BM032163.D
 Acq On : 25 Sep 2021 21:12
 Operator : CG/JU
 Sample : PB139144BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK144

Quant Time: Sep 26 01:27:33 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	7.666	152	335894	20.000	ng/u1	0.00
20) Naphthalene-d8	10.460	136	1435220	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.313	164	917342	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.036	188	1907274	20.000	ng/u1	0.00
79) Chrysene-d12	21.195	240	1889723	20.000	ng/u1	0.00
88) Perylene-d12	23.365	264	1889450	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.984	96	63971	8.042	ng/uL	0.00
4) Pyridine-d5	3.413	84	899323	39.823	ng/u1	0.00
7) Phenol-d5	6.848	99	1093246	41.003	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.990	67	660942	41.227	ng/u1	0.00
11) 2-Chlorophenol-d4	7.201	132	878233	42.427	ng/u1	0.00
15) 4-Methylphenol-d8	8.389	113	882227	42.316	ng/u1	0.00
21) Nitrobenzene-d5	8.819	128	411261	47.976	ng/u1	0.00
24) 2-Nitrophenol-d4	9.542	143	455535	50.256	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.089	165	854313	41.229	ng/u1	0.00
31) 4-Chloroaniline-d4	10.589	131	1267621	45.263	ng/u1	0.00
46) Dimethylphthalate-d6	13.713	166	2730893	45.819	ng/u1	0.00
49) Acenaphthylene-d8	14.001	160	3294233	43.503	ng/u1	0.00
54) 4-Nitrophenol-d4	14.507	143	443324	45.724	ng/u1	0.00
60) Fluorene-d10	15.301	176	2310073	45.423	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.413	200	321379	35.655	ng/u1	0.00
73) Anthracene-d10	17.136	188	3671066	46.130	ng/u1	0.00
81) Pyrene-d10	19.418	212	4126167	41.499	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.230	264	3977164	44.852	ng/u1	0.00

Target Compounds	Qvalue

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

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