

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
 Data File : BM030854.D
 Acq On : 07 Jul 2021 13:23
 Operator : CG/JU
 Sample : PB137514BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK514

Quant Time: Jul 07 13:54:47 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 06 14:32:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	89193	20.000	ng/u1	0.00
20) Naphthalene-d8	10.545	136	377033	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.392	164	233167	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.133	188	435099	20.000	ng/u1	0.00
79) Chrysene-d12	21.303	240	326506	20.000	ng/u1	0.00
88) Perylene-d12	23.550	264	346662	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	15993	7.007	ng/uL	0.00
4) Pyridine-d5	3.669	84	200504	33.590	ng/u1	0.00
7) Phenol-d5	6.934	99	275931	35.405	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.098	67	186799	37.210	ng/u1	0.00
11) 2-Chlorophenol-d4	7.298	132	230311	38.518	ng/u1	0.00
15) 4-Methylphenol-d8	8.463	113	225141	36.047	ng/u1	0.00
21) Nitrobenzene-d5	8.922	128	108831	38.810	ng/u1	0.00
24) 2-Nitrophenol-d4	9.639	143	119038	38.796	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	214355	36.303	ng/u1	0.00
31) 4-Chloroaniline-d4	10.686	131	303151	36.153	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	655620	40.163	ng/u1	0.00
49) Acenaphthylene-d8	14.086	160	821335	39.832	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	83051	29.235	ng/u1	0.00
60) Fluorene-d10	15.386	176	562493	39.133	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.515	200	69025	24.741	ng/u1	0.00
73) Anthracene-d10	17.233	188	795329	39.226	ng/u1	0.00
81) Pyrene-d10	19.521	212	768849	42.696	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.403	264	739187	41.309	ng/u1	0.00

Target Compounds Qvalue

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

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