

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083123\
 Data File : BP016890.D
 Acq On : 31 Aug 2023 10:46
 Operator : MA/JU
 Sample : PB155141BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK141

Quant Time: Sep 01 01:03:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 30 02:25:00 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.017	152	335903	20.000	ng/ul	0.00
20) Naphthalene-d8	10.846	136	1543497	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.675	164	968419	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.440	188	2067435	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1435267	20.000	ng/ul	0.00
88) Perylene-d12	24.110	264	1280845	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	69432	8.543	ng/uL	0.00
4) Pyridine-d5	3.811	84	850667	34.250	ng/ul	0.00
7) Phenol-d5	7.164	99	1017543	34.067	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	657871	34.558	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	759097	33.802	ng/ul	0.00
15) 4-Methylphenol-d8	8.722	113	813392	33.359	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	394865	33.835	ng/ul	0.00
24) 2-Nitrophenol-d4	9.928	143	418132	33.991	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.452	165	745867	32.523	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.005	131	1138114	32.361	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	2478021	33.931	ng/ul	0.00
49) Acenaphthylene-d8	14.369	160	2804056	33.964	ng/ul	0.00
54) 4-Nitrophenol-d4	14.881	143	335981	22.258	ng/ul	0.00
60) Fluorene-d10	15.669	176	2152527	35.000	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.793	200	240976	19.572	ng/ul	0.00
73) Anthracene-d10	17.540	188	3417205	37.289	ng/ul	0.00
81) Pyrene-d10	19.775	212	3627042	46.013	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.939	264	2441087	38.048	ng/ul	0.00

Target Compounds Qvalue

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

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