

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083123\
 Data File : BP016889.D
 Acq On : 31 Aug 2023 10:12
 Operator : MA/JU
 Sample : PB155062BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK062

Quant Time: Sep 01 01:01:17 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	8.016	152	376581	20.000	ng/u1	0.00
20) Naphthalene-d8	10.846	136	1696335	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.675	164	1068296	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.439	188	2320235	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	1652826	20.000	ng/u1	0.00
88) Perylene-d12	24.104	264	1504665	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	67884	7.450	ng/uL	0.00
4) Pyridine-d5	3.811	84	816095	29.309	ng/u1	0.00
7) Phenol-d5	7.164	99	977966	29.205	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	623551	29.217	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	734570	29.176	ng/u1	-0.01
15) 4-Methylphenol-d8	8.722	113	762337	27.888	ng/u1	0.00
21) Nitrobenzene-d5	9.216	128	370286	28.870	ng/u1	0.00
24) 2-Nitrophenol-d4	9.928	143	397122	29.374	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.452	165	719329	28.540	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.005	131	1063587	27.517	ng/u1	0.00
46) Dimethylphthalate-d6	14.087	166	2362898	29.329	ng/u1	0.00
49) Acenaphthylene-d8	14.375	160	2673807	29.358	ng/u1	0.00
54) 4-Nitrophenol-d4	14.881	143	327509	19.668	ng/u1	0.00
60) Fluorene-d10	15.669	176	2062874	30.406	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	228648	16.547	ng/u1	0.00
73) Anthracene-d10	17.539	188	3274636	31.840	ng/u1	0.00
81) Pyrene-d10	19.775	212	3480305	38.339	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.939	264	2379615	31.572	ng/u1	0.00

Target Compounds Qvalue

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

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