

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
 Data File : BP010916.D
 Acq On : 30 Jun 2022 12:37
 Operator : CG/JU
 Sample : N3523-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 01 02:32:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 27 23:12:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	402963	20.000	ng/u1	0.00
20) Naphthalene-d8	10.499	136	1599469	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	795734	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.122	188	1474436	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.239	240	1495523	20.000	ng/u1	# 0.00
88) Perylene-d12	23.598	264	1644041	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	50914	4.780	ng/uL	0.00
4) Pyridine-d5	3.617	84	147808	5.213	ng/u1	0.00
7) Phenol-d5	6.881	99	218025	6.835	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	642116	30.780	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	653170	24.680	ng/u1	0.00
15) 4-Methylphenol-d8	8.416	113	392955	15.583	ng/u1	-0.01
21) Nitrobenzene-d5	8.863	128	368014	36.143	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	327733	38.905	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	562516	27.025	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	613715	18.063	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	1701494	31.215	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	2140405	30.767	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	49495	5.813	ng/u1	0.00
60) Fluorene-d10	15.357	176	1520438	32.504	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.481	200	151037	30.919	ng/u1	0.00
73) Anthracene-d10	17.222	188	2534648	38.638	ng/u1	-0.01
81) Pyrene-d10	19.475	212	2898149	35.570	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.445	264	3293926	40.545	ng/u1	-0.01

Target Compounds	Qvalue

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

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