

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
 Data File : BP010915.D
 Acq On : 30 Jun 2022 12:04
 Operator : CG/JU
 Sample : N3523-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 01 02:32:29 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	415674	20.000	ng/u1	0.00
20) Naphthalene-d8	10.499	136	1623488	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	825241	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.122	188	1563658	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.239	240	1637718	20.000	ng/u1	# 0.00
88) Perylene-d12	23.598	264	1755833	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	48095	4.377	ng/uL	0.00
4) Pyridine-d5	3.617	84	196218	6.708	ng/u1	0.00
7) Phenol-d5	6.881	99	209983	6.381	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	613383	28.503	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	624077	22.860	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	372402	14.317	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	353438	34.198	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	320223	37.451	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	549834	26.025	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	563000	16.325	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	1692915	29.947	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	2064167	28.610	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	49794	5.639	ng/u1	0.00
60) Fluorene-d10	15.357	176	1515861	31.248	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.481	200	136300	26.310	ng/u1	0.00
73) Anthracene-d10	17.222	188	2382414	34.245	ng/u1	-0.01
81) Pyrene-d10	19.475	212	2764191	30.981	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.445	264	3061385	35.283	ng/u1	-0.01

Target Compounds	Qvalue

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

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