

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP011007.D
 Acq On : 18 Jul 2022 23:19
 Operator : CG/JU
 Sample : N3710-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 19 00:47:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 14 15:34:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	333299	20.000	ng/u1	0.00
20) Naphthalene-d8	10.475	136	1347621	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.340	164	730964	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.104	188	1438704	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.227	240	1451398	20.000	ng/u1	#-0.01
88) Perylene-d12	23.580	264	1488579	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	51754	5.598	ng/uL	0.00
4) Pyridine-d5	3.593	84	302660	12.665	ng/u1	0.00
7) Phenol-d5	6.858	99	210975	7.711	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	604274	34.347	ng/u1	0.00
11) 2-Chlorophenol-d4	7.210	132	605046	27.369	ng/u1	-0.01
15) 4-Methylphenol-d8	8.393	113	383697	17.518	ng/u1	-0.01
21) Nitrobenzene-d5	8.840	128	340726	34.694	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	314805	32.395	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.099	165	552610	31.095	ng/u1	0.00
31) 4-Chloroaniline-d4	10.622	131	595257	20.317	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	2023550	40.504	ng/u1	0.00
49) Acenaphthylene-d8	14.034	160	2243240	38.056	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	55337	5.473	ng/u1	0.00
60) Fluorene-d10	15.339	176	1691185	39.825	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200	207656	29.169	ng/u1	-0.01
73) Anthracene-d10	17.204	188	2642527	43.205	ng/u1	-0.01
81) Pyrene-d10	19.463	212	3041283	39.897	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.427	264	3116093	43.585	ng/u1	-0.02

Target Compounds	Qvalue

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

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