

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
 Data File : BP010807.D
 Acq On : 24 Jun 2022 22:55
 Operator : CG/JU
 Sample : N3401-15
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AG1

Quant Time: Jun 24 23:41:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	340667	20.000	ng/u1	0.00
20) Naphthalene-d8	10.516	136	1391867	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	783807	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.139	188	1567746	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.251	240	1556058	20.000	ng/u1	# 0.00
88) Perylene-d12	23.621	264	1615386	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	48881	5.428	ng/uL	0.00
4) Pyridine-d5	3.610	84	289241	12.066	ng/u1	0.00
7) Phenol-d5	6.887	99	234885	8.710	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	693870	39.343	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	691485	30.906	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	437646	20.529	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	412288	46.531	ng/u1	0.00
24) 2-Nitrophenol-d4	9.604	143	361016	49.248	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	666210	36.781	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	998483	33.771	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	2328953	43.377	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2850942	41.604	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	62183	7.415	ng/u1	0.00
60) Fluorene-d10	15.375	176	1981919	43.015	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	176309	33.944	ng/u1	0.00
73) Anthracene-d10	17.239	188	3055661	43.808	ng/u1	0.00
81) Pyrene-d10	19.486	212	3484357	41.101	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.468	264	3557221	44.563	ng/u1	0.00

Target Compounds Qvalue

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

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