

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010875.D
 Acq On : 27 Jun 2022 13:42
 Operator : CG/JU
 Sample : N3392-14 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jun 27 23:15:22 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	7.710	152	390597	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	1538933	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	771499	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.133	188	1383280	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.245	240	1366406	20.000	ng/u1	# 0.00
88) Perylene-d12	23.610	264	1544787	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	30501	2.954	ng/uL	0.00
4) Pyridine-d5	3.611	84	360578	13.119	ng/u1	0.00
7) Phenol-d5	6.887	99	255295	8.257	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	761148	37.641	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	762447	29.721	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	462654	18.928	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	434401	44.341	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	387078	47.757	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	671760	33.543	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	1037074	31.724	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2108554	39.898	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	2506422	37.160	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	58368	7.071	ng/u1	0.00
60) Fluorene-d10	15.363	176	1871570	41.268	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	178637	38.979	ng/u1	0.00
73) Anthracene-d10	17.228	188	2888333	46.931	ng/u1	0.00
81) Pyrene-d10	19.480	212	3197911	42.958	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.457	264	3669746	48.073	ng/u1	0.00

Target Compounds Qvalue

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

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