

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010846.D
 Acq On : 26 Jun 2022 01:45
 Operator : CG/JU
 Sample : N3392-09 10X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY4

Quant Time: Jun 26 22:55:42 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	7.716	152	478280	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1925437	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	1003630	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1771332	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1724391	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	1866876	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	31238	2.471	ng/uL	0.00
4) Pyridine-d5	3.617	84	267664	7.953	ng/u1	0.01
7) Phenol-d5	6.893	99	273031	7.211	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	855949	34.569	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	853648	27.176	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	518775	17.333	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	517338	42.207	ng/u1	0.00
24) 2-Nitrophenol-d4	9.599	143	469627	46.311	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	797871	31.843	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	1151491	28.153	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2597932	37.789	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	3188091	36.334	ng/u1	0.00
54) 4-Nitrophenol-d4	14.587	143	60057	5.593	ng/u1	0.00
60) Fluorene-d10	15.369	176	2201737	37.319	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	174631	29.757	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3194613	40.536	ng/u1	-0.01
81) Pyrene-d10	19.480	212	3503508	37.293	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.457	264	3845613	41.686	ng/u1	-0.02

Target Compounds Qvalue

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

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