

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010840.D
 Acq On : 25 Jun 2022 22:22
 Operator : CG/JU
 Sample : N3392-06 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/27/2022
 Supervised By :mohammad ahmed 07/01/2022

Quant Time: Jun 26 22:54:42 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	497424	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1981858	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	1035380	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1898540	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1842585	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	1987834	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	25992m	1.977	ng/uL	0.00
4) Pyridine-d5	3.617	84	229938	6.569	ng/u1	0.01
7) Phenol-d5	6.893	99	234125	5.946	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	708556	27.515	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	712016	21.795	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	428969	13.781	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	425220	33.704	ng/u1	0.00
24) 2-Nitrophenol-d4	9.599	143	382646	36.659	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	649104	25.168	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	964290	22.905	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2198710	31.001	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	2566099	28.349	ng/u1	0.00
54) 4-Nitrophenol-d4	14.587	143	58313	5.264	ng/u1	0.00
60) Fluorene-d10	15.369	176	1832527	30.109	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.492	200	155261	24.684	ng/u1	0.00
73) Anthracene-d10	17.233	188	2888495	34.196	ng/u1	-0.01
81) Pyrene-d10	19.486	212	3209140	31.968	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.457	264	3465160	35.276	ng/u1	-0.02

Target Compounds Qvalue

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

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